



In response to the Office Action of May 20, 2002, please amend the application as follows:

IN THE SPECIFICATION

Between the title and "Field of Invention", insert the following:

--Cross Reference to Related Applications

This application is a national stage entry under 35 U.S.C. §371 of the PCT/EP99/05459 filed July 30, 1994.--

Page 1, lines 22-27:

B1
*should
use Ar*
B2
*should
be a
common*
-(CH₂)_r Ar₁ where r is 0, 1 or 2 and Ar is an aromatic group chosen among benzene, naphthalene, thiophene, benzothiophene, pyridine, quinoline, indole, furan, benzofuran, thiazole, benzothiazole, imidazole, benzoimidazole, possibly substituted with up to 2 substituents chosen among C₁₋₃ alkyl, C₁₋₃ haloalkyl, C₁₋₃ alkyloxy and C₂₋₄ amino-alkyloxy, halogens, OH, NH₂, CN, NR₆R₇, where R₆ and R₇, are the same or different, and are H or C₁₋₃ alkyl.

Page 1, line 29-32 to Page 1a, line 2:

*Sub
C4 B3*
-(CH₂)_rAr₁ where r is 0, 1 or 2 and Ar₁ is an aromatic group chosen among: benzene, naphthalene, thiophene, benzothiophene, pyridine, quinoline, indole, furan, benzofuran, thiazole, benzothiazole, imidazole, benzoimidazole, possibly substituted with up to 2 groups chosen among: C₁₋₃ alkyl, C₁₋₃haloalkyl, C₁₋₃ alkyloxy and C₂₋₄ amino-alkyloxy, halogens, OH, NH₂, CN, NR₆R₇, where R₆ and R₇, are the same or different and are H or C₁₋₃ alkyl.

Page 2, lines 10-12

*B4
can't be
both N's*
or R₈ and R₉ together with the N atom to which they are linked to form a piperazine possibly substituted at one of its nitrogen atoms by C₁₋₃ alkyl, C₁₋₃ acyl or methanesulfonyl;

B5
Page 3, lines 10-14

In WO9834949 it is described how compounds having lower molecular weight, monocyclic, containing only four bi-functional residues linked among each other by a peptide or pseudopeptide bond present pharmacological activity similar or higher than that of known compounds and moreover show a high selectivity for the human NK2 receptor.

Page 5, line 5 delete "tryptophane" insert --tryptophane--.

More preferred are the compounds of formula (1) wherein:

-X₁, X₂, X₃, X₄ are -CONR-,

-R is H;

-R₁ is the lateral chain of tryptophane;

-R₂ is the lateral chain of phenylalanine substituted with up to two residues selected from the group consisting of: chlorine, fluorine, CF₃, OH, and CN; or a group 3-pyridyl-methyl, or 4-pyridyl-methyl;

-R₃ is benzyl,

and the other substituents are as above defined.

B7
Sub CS
Page 5, lines 15-20

R₉ is a group chosen among: 4-tetrahydropyranyl, 4-tetraiodrothiopyranyl, 1-oxotetraiodrothiopyran-4-yl, 1,1 dioxo-tetrahydrothiopyran-4-yl, N-methyl-4-piperidinyl, N-methansulfonyl-4-piperidinyl, N-aminosulfonyl-4-piperidinyl, or R₈ and R₉ together with the N atom to which they are linked represent N-methyl-piperazinyl, N-acetyl-piperazinyl, piperazinyl, N-methanesulfonyl-piperazinyl.

B8
Page 10, line 32 and page 11, line 1

As starting compound the cyclo {Succinic[1-(R)-amino]-Trp-Phe-[(R)-NH-CH(CH₂C₆H₅)-CH₂-NH]-} (Compound A).